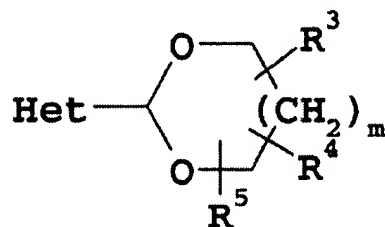


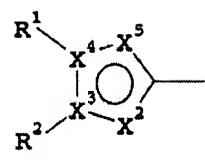
CLAIMS

1. (currently amended) A compound of formula (I):

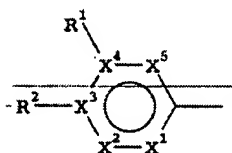


(I)

wherein:-



Het is a five ~~or six~~ membered heteroaromatic ring of the formula



in which one of R¹ is optionally substituted aryl, and R² is 4-pyridyl;

optionally substituted heteroaryl and the other is optionally substituted heteroaryl or optionally substituted aryl; wherein heteroaryl is selected from: optionally substituted benzimidazolyl, furyl, imidazolyl, isoxazolyl, isoquinolinyl, isothiazolyl, oxadiazolyl, pyrazinyl, pyridazinyl, pyrazolyl, pyridyl, pyrimidinyl, pyrrolyl, quinazolinyl, quinolinyl, 1,3,4-thiadiazolyl, thiazolyl, thienyl and triazolyl groups; and heteroaryl optional substitution is with one or more substituents selected from: acyl, acylamino, alkoxycarbonyl, alkylenedioxy, aroyl, aroylamino, aryl, arylalkyloxycarbonyl, aryloxycarbonyl, carboxy, cyano, halo, heteroaroyl, heteroaryl, heteroaroylamino, hydroxy, nitro, trifluoromethyl, R¹¹Z², Y¹Y²N, Y¹Y²N-CO, Y¹Y²NSO₂, alkylSO₂-Y¹N or alkyl optionally substituted with aryl, heteroaryl, hydroxy, oxo, CO₂R⁷, CONY³Y⁴ or NY¹Y²; wherein aryl is selected from: phenyl and naphthyl; and aryl optional substitution is with one or more substituents selected from: acyl, acylamino, alkoxy, alkoxycarbonyl, alkylenedioxy, alkylsulphinyl, alkylsulphonyl, alkylthio, aroyl, aroylamino, aryl, arylalkyloxy, arylalkyloxycarbonyl, arylalkylthio, aryloxy, aryloxycarbonyl, arylsulphinyl, arylsulphonyl, arylthio, carboxy, cyano, halo, heteroaroyl, heteroaryl, heteroarylalkyloxy,

heteroaroylamino, heteroaryloxy, hydroxy, nitro, trifluoromethyl, Y^3Y^4N- , Y^3Y^4NCO- , $Y^3Y^4NSO_2-$, $Y^3Y^4N-C_{2-6}alkylene-Z^1-$ (where Z^1 is O, NR^5 or $S(O)_n$), alkylC(=O)- Y^3N- , alkylSO₂- Y^3N- or alkyl optionally substituted with aryl, heteroaryl, hydroxy, or Y^3Y^4N- ; X^1 is a bond, X^3 and X^4 are each independently N or C and X^2 and X^5 are independently CH, N, NH, O or S; or X^3 and X^4 are C, one of X^1 , X^2 and X^5 is N and the others are N or CH; but excluding compounds in which X^1 is a bond, one of X^2 and X^5 is N and the other is NH and X^3 and X^4 are both C X^1 is CH, X^3 is C, X^4 is N and X^5 is N;

R^3 represents a group -L¹- R^6 ;

R^4 represents hydrogen, alkyl or hydroxyalkyl; or

R^3 and R^4 , when attached to the same carbon atom, may form with the said carbon atom a cycloalkyl, cycloalkenyl or heterocycloalkyl ring or a group C=CH₂;

R^5 represents hydrogen or alkyl;

R^6 is hydrogen, alkyl, azido, hydroxy, alkoxy, aryl, arylalkyloxy, aryloxy, carboxy, an acid bioisostere selected from the group consisting of C(=O)NHOH, -C(=O)-CH₂OH, -C(=O)-CH₂SH, C(=O)NH-CN, sulpho, phosphono, alkylsulphonylcarbamoyl, tetrazolyl, arylsulphonylcarbamoyl, heteroarylsulphonylcarbamoyl, N methoxycarbamoyl, 3 hydroxy-3-cyclobutene-1,2-dione, 3,5-dioxo-1,2,4-oxadiazolidinyl, 3 hydroxyisoxazolyl and 3 hydroxy 1 methylpyrazolyl, cycloalkyl, cycloalkyloxy, heteroaryl, heteroarylalkyloxy, heteroaryloxy, heterocycloalkyl, heterocycloalkyloxy, nitro, -NY¹Y², -N(R^7)-C(=Z)- R^8 , -N(R^7)-C(=Z)-L²- R^9 , -NH-C(=Z)-NH- R^8 , -NH-C(=Z)-NH-L²- R^9 , -N(R^7)-SO₂- R^8 , -N(R^7)-SO₂-L²- R^9 , -S(O)_n R^{10} , -C(=Z)-NY¹Y² or -C(=Z)-OR¹⁰;

R^7 is hydrogen, alkyl, aryl, arylalkyl, cycloalkyl, heteroaryl, heteroarylalkyl, or heterocycloalkyl;

R^8 is alkyl, alkoxy, aryl, arylalkyloxy, cycloalkyl, heteroaryl, heteroarylalkyloxy or heterocycloalkyl;

R^9 is alkoxy, aryl, arylalkyloxy, arylalkyloxycarbonylamino, carboxy, an acid bioisostere selected from the group consisting of C(=O)NHOH, -C(=O)-CH₂OH, -C(=O)-CH₂SH, C(=O)NH-CN, sulpho, phosphono, alkylsulphonylcarbamoyl, tetrazolyl, arylsulphonylcarbamoyl, heteroarylsulphonylcarbamoyl, N methoxycarbamoyl, 3 hydroxy-3-cyclobutene-1,2-dione, 3,5-dioxo-1,2,4-oxadiazolidinyl, 3 hydroxyisoxazolyl and 3 hydroxy 1 methylpyrazolyl, cycloalkyl, cyano, halo, heteroaryl, heteroarylalkoxy, heterocycloalkyl, hydroxy or -NY³Y⁴;

R^{10} is alkyl, aryl, arylalkyl, cycloalkyl, heteroaryl, heteroarylalkyl, or heterocycloalkyl;

L^1 represents a direct bond or a straight- or branched-chain alkylene linkage containing from 1 to 6 carbon atoms and optionally substituted by halogen, hydroxy, alkoxy or oxo;

L^2 is a straight- or branched-chain alkylene linkage containing from 1 to 6 carbon atoms;

Y^1 and Y^2 are independently hydrogen, alkenyl, alkynyl, aryl, cycloalkyl, heterocycloalkyl, heteroaryl or alkyl optionally substituted by alkoxy, aryl, cyano, cycloalkyl, heteroaryl, heterocycloalkyl, hydroxy, oxo, $-CO_2R^7$, $-CONY^3Y^4$ or $-NY^3Y^4$, or the group $-NY^1Y^2$ may form a 5-7 membered cyclic amine which (i) may be optionally substituted with one or more substituents selected from alkoxy, carboxamido, carboxy, hydroxy, oxo (or a 5, 6, or 7 membered cyclic acetal derivative thereof), alkyl, aryl, arylalkyl, cycloalkyl, heteroaryl, heteroarylalkyl, or heterocycloalkyl or alkyl substituted by carboxy, carboxamido or hydroxy (ii) may also contain a further heteroatom selected from O, S, SO_2 or NY^5 and (iii) may also be fused to additional aryl, heteroaryl, heterocycloalkyl or cycloalkyl rings to form a bicyclic or tricyclic ring system; Y^3 and Y^4 are independently hydrogen, alkenyl, alkyl, alkynyl, aryl, arylalkyl, cycloalkyl, heteroaryl or heteroarylalkyl, or the group $-NY^3Y^4$ may form a 5-7 membered cyclic amine as defined for $-NY^1Y^2$ above;

Y^5 is hydrogen, alkyl, aryl, arylalkyl, $-C(=Z)R^{10}$, $-C(=Z)OR^{10}$ or $-SO_2R^{10}$;

Z is an oxygen or sulphur atom;

m is zero or an integer 1 or 2; and

n is zero or an integer 1 or 2;

~~or and an N-oxide thereof, or and an ester prodrug thereof, or and a pharmaceutically acceptable salt thereof, or and a hydrate of a compound of formula (I), or and an N-oxide thereof, and its ester prodrug.~~

2. (cancelled)

3. (cancelled)

4. (cancelled)

5. (previously presented) A compound according to Claim 1 in which one of R^1 and R^2 is 4-pyridyl and the other is 4-fluorophenyl.

6. (cancelled)

7. (cancelled)

8. (cancelled)

9. (cancelled)

10. (cancelled)

11. (previously presented) A compound according to Claim 1 in which R³ and R⁴ are both C₁₋₄alkyl groups.

12. (previously presented) A compound according to Claim 1 in which R³ is -C(=O)-NY¹Y² (where Y¹ and Y² are as defined in Claim 1) and R⁴ is C₁₋₄alkyl.

13. (previously presented) A compound according to Claim 12 in which Y¹ is hydrogen and Y² is alkyl or cycloalkyl.

14. (cancelled)

15. (previously presented) A pharmaceutical composition comprising a compound according to Claim 1 together with a pharmaceutically acceptable carrier or excipient.

16-20 (cancelled)